

# **Fuels Workshop on Regulatory and Non-Regulatory Fuels Activities for 2006**

**September 22, 2006**

**California Environmental Protection Agency**

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**Air Resources Board**

# Agenda

- ➡ Introductions and Schedule
- ➡ California Predictive Model
  - Reactivity of Evaporative and Exhaust Emissions
    - CO Reactivity
  - 2006 Draft Predictive Model
  - Emission Inventory
- ➡ Presentations by Others
- ➡ Open Discussions
- ➡ Closing Remarks

# Tentative Future 2006 Workshops

- ➡ October 6    9:00 to 12:30    Sher Auditorium
- ➡ October 25    9:00 to 12:30    Sher Auditorium

- The October workshops will be webcast. All meetings will be available by conference call.
- Next set of workshops will be scheduled based on progress

# Reactivity of Evaporative and Exhaust Emissions

# Reactivity Progress

## 2006 Update to Draft MIR Values

- ➡ 2006 MIR list was presented to the Reactivity Subgroup and used in calculating reactivities
- ➡ Data sets for Diurnal, Hot Soak and Exhaust obtained from in-use testing at El Monte
- ➡ All speciated data was for E6 fuel
- ➡ Running loss was calculated from headspace and liquid fuel profiles (per Dr. Harley, UCB)
- ➡ Presented to Reactivity Working Group for review

# 2006 Draft Specific Reactivity

(applied to speciated data sets from VEDS database)

	2006 Draft Specific Reactivity
Exhaust	3.99
Hot Soak	3.12
Diurnal	2.36
Running Loss	2.54

# Updates to Predictive Model Related MIR Information

<http://www.arb.ca.gov/fuels/gasoline/premodel/pmdevelop.htm#MIR>

# CO Reactivity



## Reactivity Values for Predictive Model

- ➡ Maximum Incremental Reactivity (MIR) values are used in the Predictive Model (PM)
- ➡ Consistent with the previous PM assessment
- ➡ First developed in early 1990 by Dr. Carter at UCR and updated several times since then
- ➡ MIR is deemed most appropriate for scientific and regulatory applications by the Reactivity Research Scientific Committee

## MIR (continued)

- ➡ The Tables of MIR Values were adopted by ARB in June 2000 and updated in December 2003 to ensure that our regulation is based on the best sound science.
- ➡ Required to review the Tables of MIR Values every 18 months to determine if modifications to the values are warranted.
- ➡ Used in Low Emission Vehicle and Clean Fuel (LEV/CF) and aerosol coatings regulations, and possibly for other categories

## MIR (Continued)

- ➡ 3-D airshed model derived reactivity values would be the most appropriate but are available only for a limited number of VOCs (~30) due to computational demands
- ➡ Comparison study between 3-D and box model derived reactivity values indicates that the correlation is high
- ➡ The MIR-based California aerosol coatings regulation was approved by the U.S. EPA in January 2005

## MIR (Continued)

- ➡ U.S. EPA publishes an interim guideline in September 2005 on VOC reactivity and encourages all states to consider it in development of ozone SIPs
- ➡ Other agencies are using MIR scale as VOC control strategies
- ➡ Working with Reactivity Research Working Group to develop other reactivity metrics
- ➡ MIR is the most scientifically sound reactivity scale available for ~800 VOCs

## CO Reactivity

- ☞ Treated as a VOC in SAPRC99 mechanism and listed in the Tables of MIR Values
- ☞ Is a slow reacting chemical so the box model derived MIR value for CO may be an underestimation
- ☞ Included in the comparison study and its relative reactivity is consistent in terms of rankings
- ☞ Inappropriate to use different reactivity scales for any reactivity applications (MIR-3D for CO vs. MIR for others)
- ☞ The MIR value (0.06) for CO is appropriate for the predictive model.

# 2006 Draft Predictive Model

# 2006 Draft Predictive Model

The 2006 Draft Predictive Model includes several major revisions:

- ☞ Draft statistical models for exhaust THC, NO<sub>x</sub> and CO.
- ☞ The 2010 vehicle emission weights from the EMFAC 2007 working draft model, including permeation estimates, using California 8-hour temperature profile and relative humidity.
- ☞ Updated Maximum Incremental Reactivity (MIR) values :
  - Based on the 2006 list of MIR of total organic gaseous compounds.
  - Used to calculate reactivities of exhaust and evaporative processes
  - Ethanol permeation reactivity was based on the CRC E-65 study.

# Predictive Model Fundamental Eqn

$$\%ChangeinMassEmission = \frac{(Emission_{Cand} - Emission_{Ref})}{Emission_{Ref}} \times 100\%$$

- ➡ This fundamental equation has never changed since the Predictive Model adopted by the Board
- ➡ The equation also applies to permeation
- ➡ The MIR is used to provide flexibility for refiners to offset exhaust hydrocarbon emissions with evaporative hydrocarbon emissions.

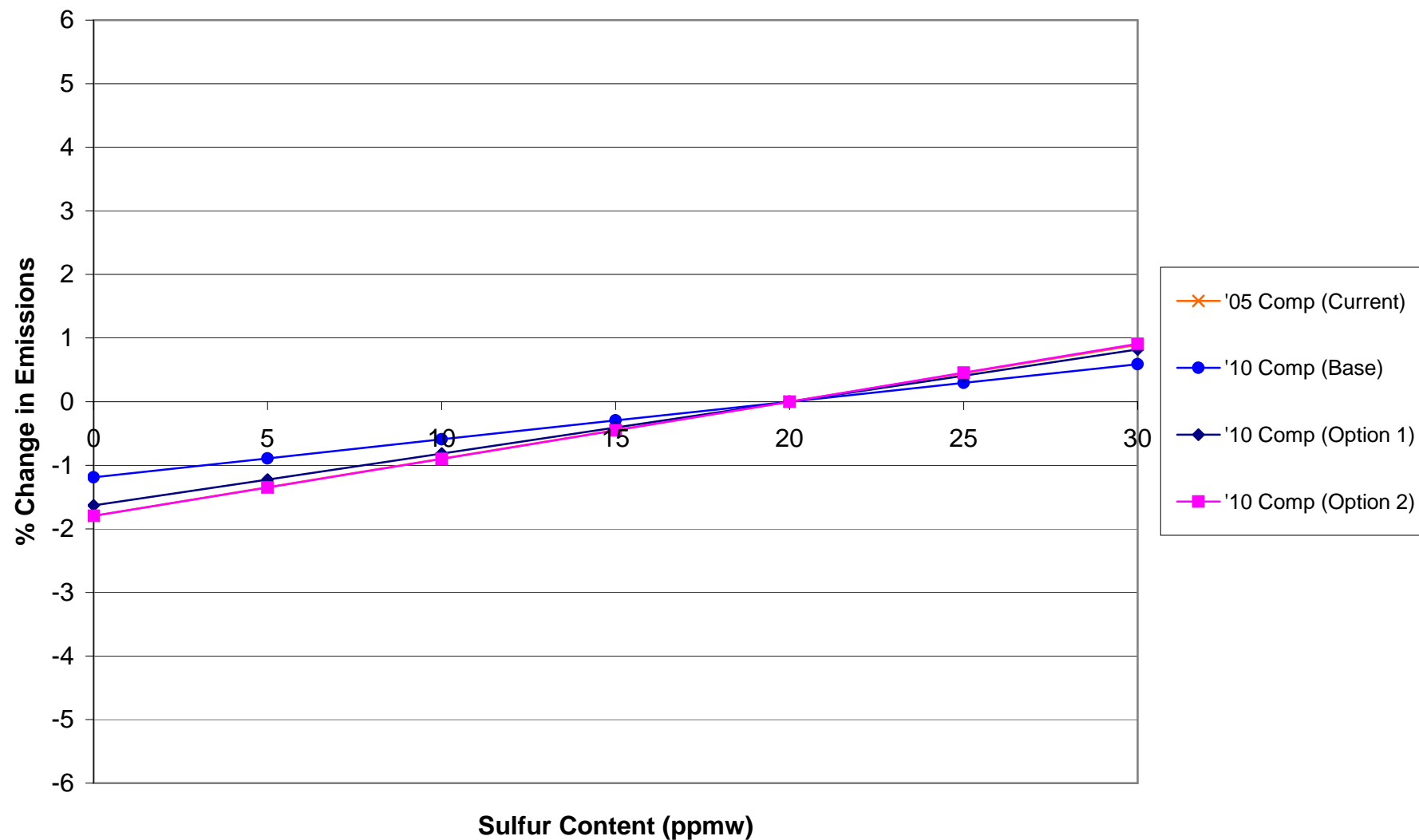


## 2006 Draft Predictive Model (2010 Base Year)

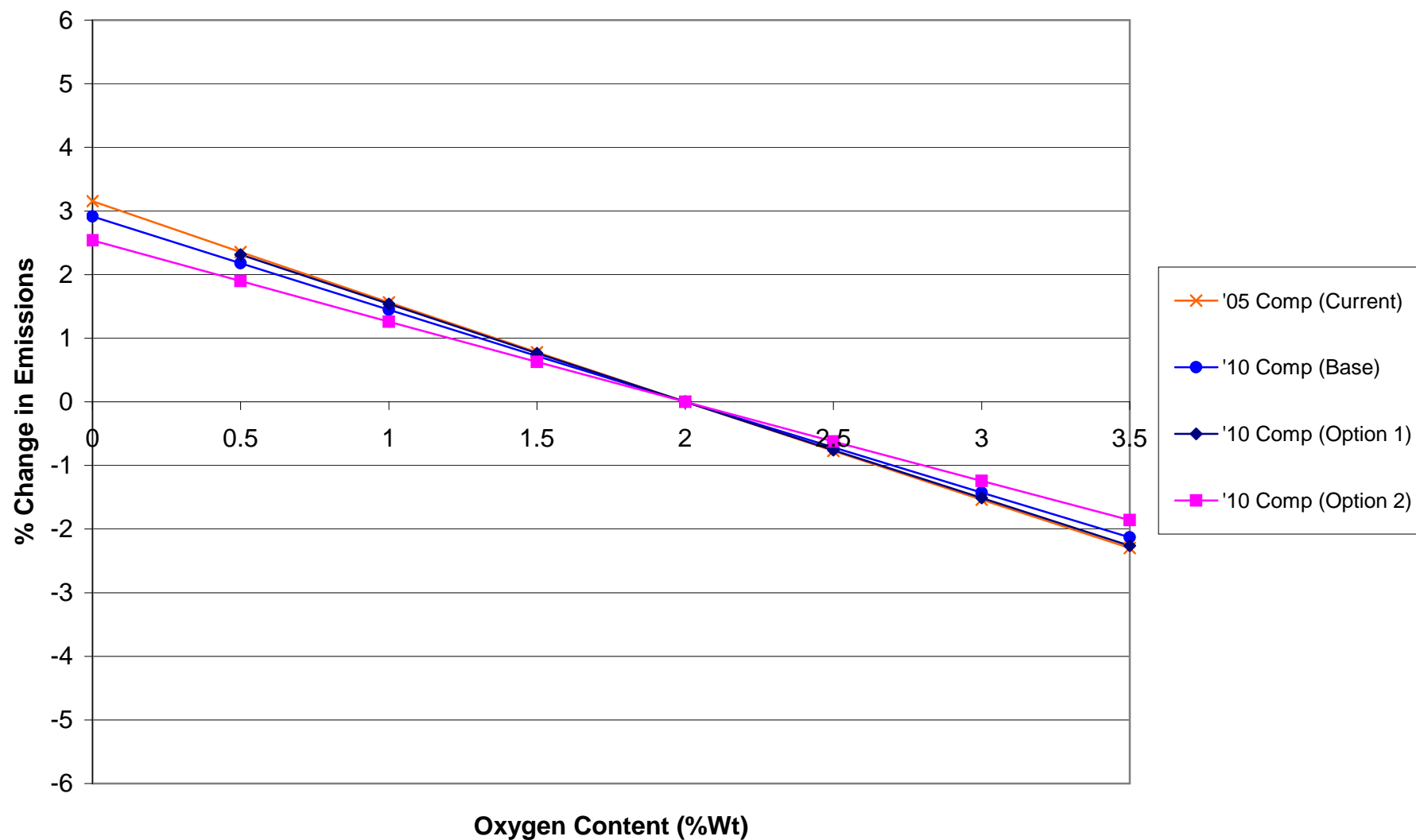
Pollutant	Emission (tpd)	MIR	OFP	
			(tpd)	(%)
Exh TOG	249	3.99	994	43.6
CO	4378	0.06	263	11.5
Evap TOG				
DI/RT	118	2.36	278	12.2
HS	64	3.12	200	8.8
RL	170	2.76	469	20.6
Perm	23	3.27	75	3.3

# THC Response to Fuel Properties

### THC Response to Sulfur (All Other Fuel Properties @ Flat Limits)

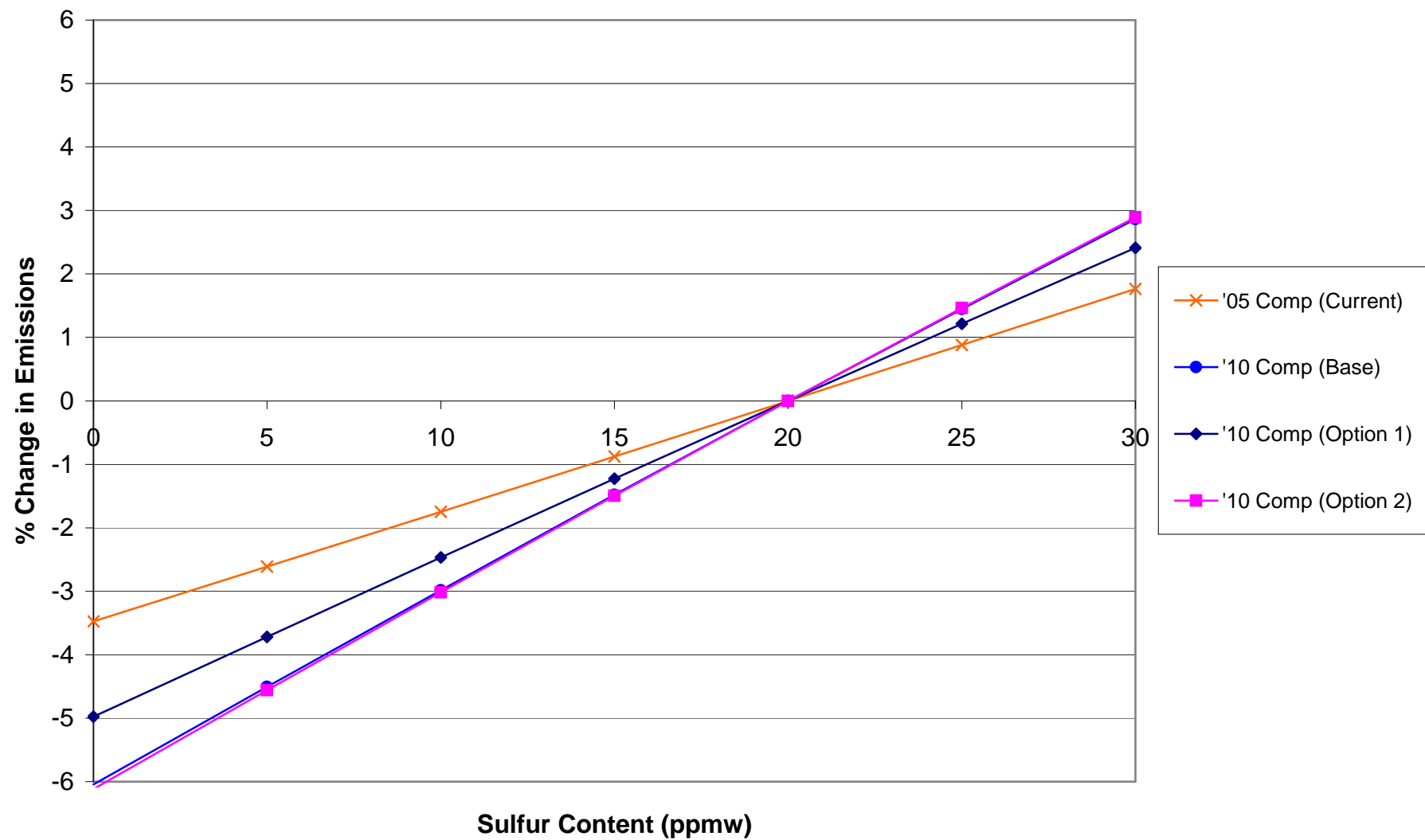


### THC Response to Oxygen (All Other Fuel Properties @ Flat Limits)

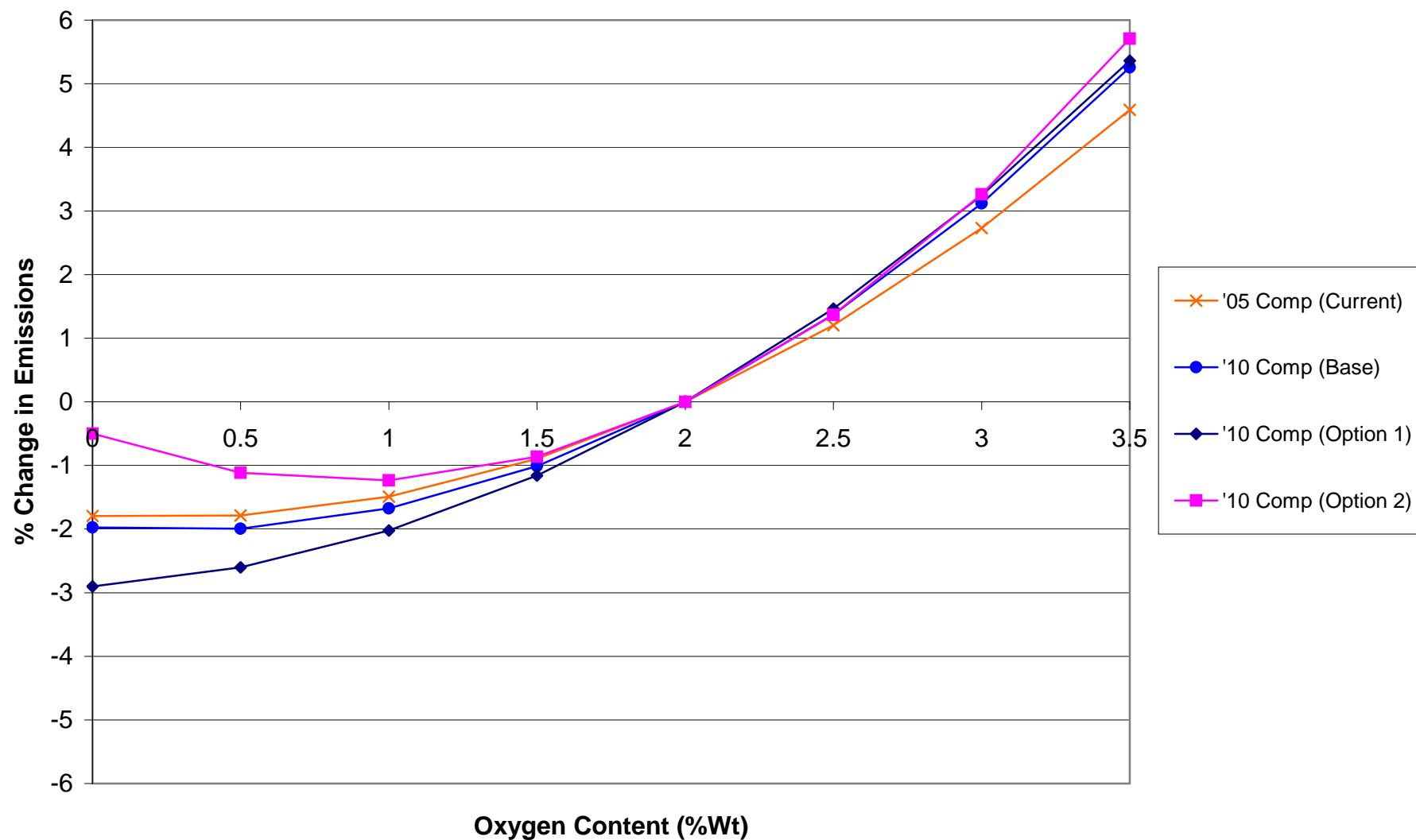


# **NO<sub>x</sub> Response to Fuel Properties**

### NOx Response to Sulfur (All Other Fuel Properties @ Flat Limits)



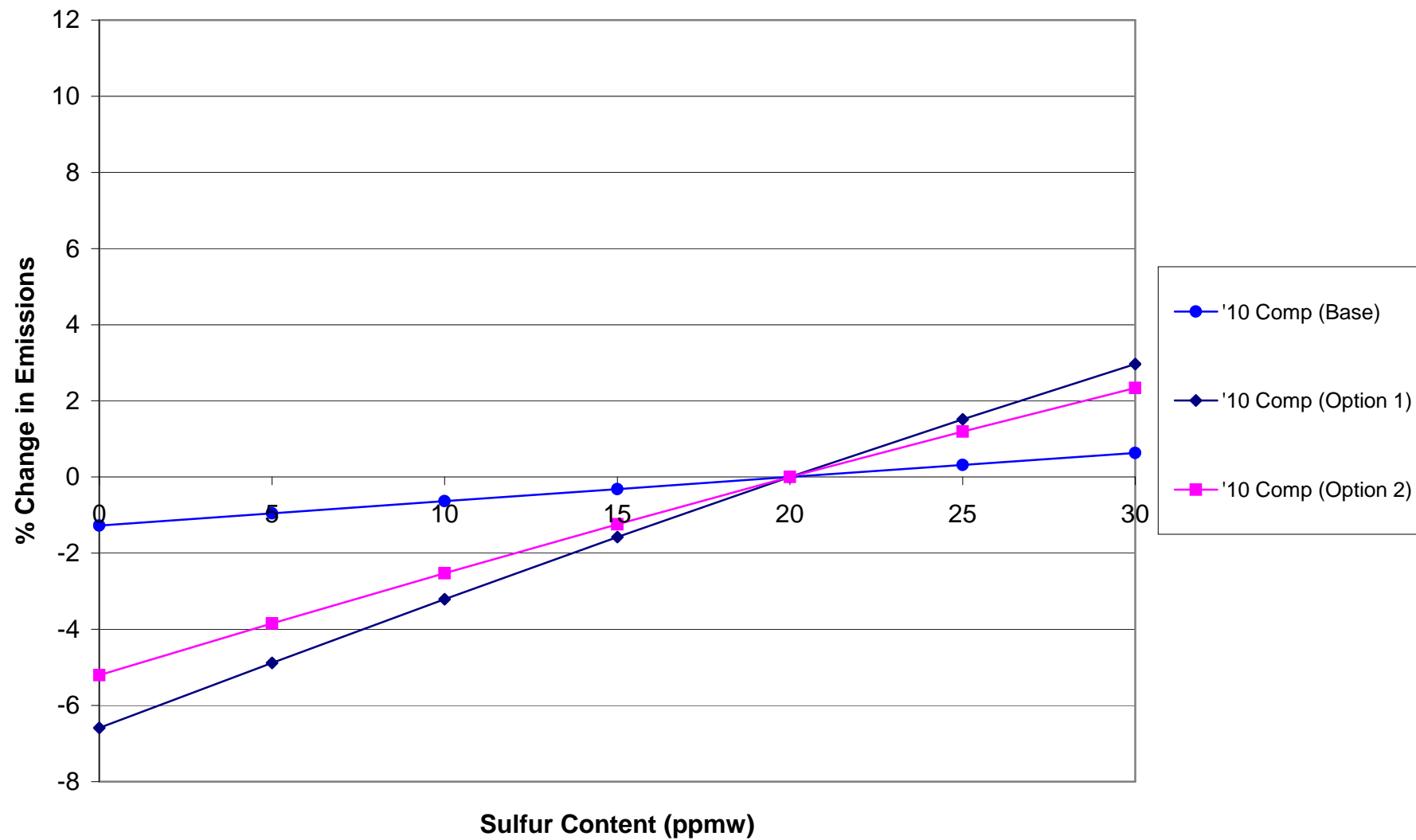
### NOx Response to Oxygen (All Other Fuel Properties @ Flat Limits)



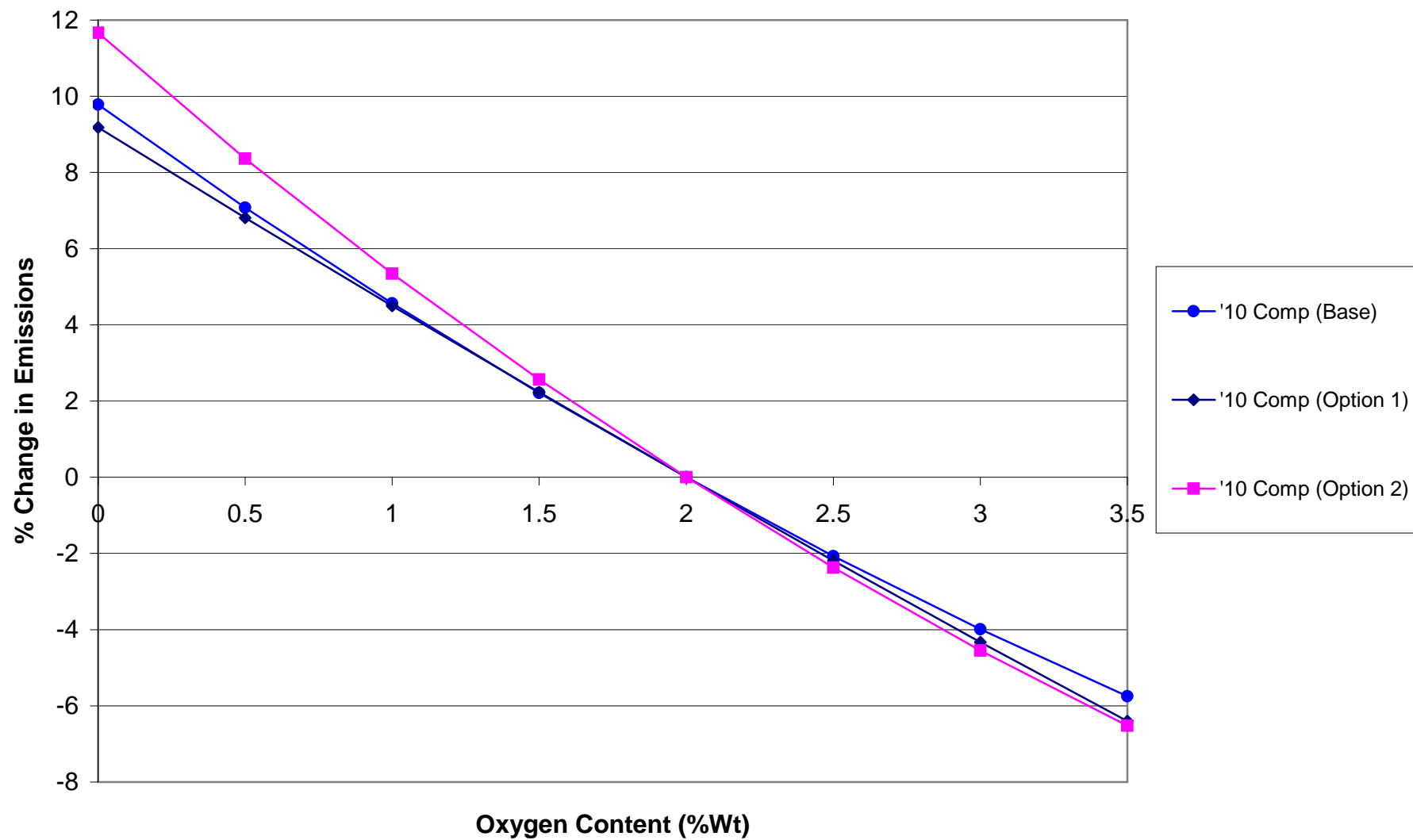
# CO Response to Fuel Properties



### CO Response to Sulfur (All Other Fuel Properties @ Flat Limits)



### CO Response to Oxygen (All Other Fuel Properties @ Flat Limits)



# Emissions Inventory

# **Presentations by Others**

# Open Discussions

# Closing Remarks